

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062515\
 Data File : BF080183.D
 Acq On : 26 Jun 2015 00:55
 Operator : TP/IZ
 Sample : G2699-01RE
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LAR9819RE

Manual Integrations
 APPROVED

apatel
 6/26/2015 2:06:01 PM

Quant Time: Jun 26 03:49:19 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 26 01:01:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.29	152	29092	20.00	ng	0.00
21) Naphthalene-d8	8.58	136	115939	20.00	ng	0.00
38) Acenaphthene-d10	10.36	164	64320	20.00	ng	0.00
63) Phenanthrene-d10	11.86	188	128862	20.00	ng	0.00
75) Chrysene-d12	14.77	240	136347	20.00	ng	-0.06
86) Perylene-d12	16.62	264	133337	20.00	ng	-0.09
System Monitoring Compounds						
5) 2-Fluorophenol	5.90	112	264922	161.20	ng	0.00
7) Phenol-d6	6.89	99	342294	156.52	ng	-0.01
23) Nitrobenzene-d5	7.85	82	218840	109.89	ng	0.00
41) 2,4,6-Tribromophenol	11.16	330	104799	166.63	ng	0.00
44) 2-Fluorobiphenyl	9.66	172	401411	89.00	ng	-0.01
78) Terphenyl-d14	13.52	244	499838	85.62	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.61	128	13408	2.21	ng	98
49) Dimethylphthalate	10.05	163	43743	9.05	ng	# 97
70) Phenanthrene	11.89	178	17387	2.23	ng	97
74) Fluoranthene	13.13	202	44490	5.35	ng	87
77) Pyrene	13.38	202	43809	5.22	ng	97
80) Benzo(a)anthracene	14.76	228	38908	4.68	ng	94
82) Chrysene	14.80	228	32491	4.24	ng	95
85) Indeno(1,2,3-cd)pyrene	18.44	276	34574	3.58	ng	# 84
87) Benzo(b)fluoranthene	16.08	252	57235m	7.09	ng	
88) Benzo(k)fluoranthene	16.12	252	17166m	2.09	ng	
89) Benzo(a)pyrene	16.54	252	44651	5.82	ng	# 85
91) Benzo(g,h,i)perylene	18.99	276	34434	4.35	ng	# 78

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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